Application No.: 10/583,469

## AMENDMENTS TO THE CLAIMS

This listing of claims will replace all prior versions and listings of claims in the application:

#### LISTING OF CLAIMS:

1. (Currently Amended) A compound of formula (I)

wherein ring A represents a benzene ring which may have a substituent(s), wherein the substituent(s) is 1 to 5 optional substituent(s) selected from the group of methyl, ethyl, a fluorine atom, a chlorine atom, methoxy, ethoxy, diffuoromethoxy, hydroxyl, and acetyl, trifluoromethoxy, methylsulfonyl, acetylamino, methylsulfonylamino, 1 hydroxy 1 methylethyl, 1 propenyl and eyano;

ring B represents a benzene ring which may have a substituent(s), a thiophene ring which may have a substituent(s), or a 1,3-benzodioxole ring which may have a substituent(s), or a cyclopentane ring which may have a substituent(s), wherein the substituent(s) is 1 to 5 optional substituent(s) selected from the group consisting of methyl, ethyl, propyl, a fluorine atom, and a chlorine atom, methoxy and ethoxy;

K represents C1-4 alkylene which may be substituted with 1 to 5 optional substituent(s) selected from the group consisting of methyl, a fluorine atom, hydroxy, and oxo;

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O represents methylene, or ethylene, -O-, or -CH2-O-;

M represents a bond, C1-4 alkylene which may have a substituent(s), or C2-4 alkenylene which may have a substituent(s), wherein the substituent(s) is 1 to 5 optional substituent(s) selected from the group consisting of methyl-and hydroxyl;

ring D represents a benzene ring which may have a substituent(s), a pyrrole ring which may have a substituent(s), an imidazole ring which may have a substituent(s), an oxazole ring which may have a substituent(s), or a thiophene ring which may have a substituent(s), or a thiophene ring which may have a substituent(s), wherein the substituent(s) is 1 to 5 optional substituent(s) selected from the group consisting of earboxy, methyl, a fluorine atom, a chlorine atom, and methoxycarbonyl, ethoxycarbonyl, aminocarbonyl, N-methylaminocarbonyl, NN-dimethylaminocarbonyl, and acetyl;

ring E represents a benzene ring-which may have a substituent(s), wherein the substituent(s) is 1 to 5 optional substituent(s) selected from the group consisting of methyl, a chlorine atom, a fluorine atom, methoxy, and ethoxy;

L represents a bond, or CH2, O, S, SO, SO2 or NH-;

Z represents COOH, CONHSO<sub>2</sub>R<sup>1</sup>, in which R<sup>1</sup> represents <u>methyl, or C1-8 alkyl which</u> may be substituted, a benzene ring which may have a substituent(s), a pyridine ring which may have a substituent(s), a furan ring which may have a substituent(s), a furan ring which may have a substituent(s), a thiazole ring which may have a substituent(s), a morpholine ring which may have a substituent(s), a morpholine ring which may have a substituent(s), a morpholine ring which may have a substituent(s), wherein the substituent(s) is 1 to 5 optional substituent(s)

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selected from the group consisting of methyl, tert-butyl, a chlorine atom, a fluorine atom, trifluoromethyl, methoxy, trifluoromethoxy, and acetyl; and

t represents 0 or 1, or

a salt thereof, or a solvate thereof.

(Original) The compound according to claim 1, wherein the compound of formula (I) is an optically active compound of formula (I-A):

wherein represents  $\beta$ -configuration; and other symbols have the same meanings as described in claim 1.

## 3.-4. (Canceled)

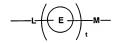
(Previously presented) The compound according to claim 1, wherein ring B is an indan ring which may have a substituent(s).

# 6. (Canceled)

7. (Original) The compound according to claim 1, wherein ring D is a benzene ring which may have a substituent(s), a pyrazole ring which may have a substituent(s) or a pyrrole ring which may have a substituent(s).

#### 8. (Canceled)

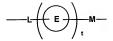
9. (Original) The compound according to claim 1, wherein



is methylene which may be substituted, ethylene which may be substituted, propylene which may be substituted, or ethenylene which may be substituted.

10. (Currently Amended) The compound according to claim 1, wherein ring A is a benzene ring which may have a substituent(s); ring B is an indan ring which may have a substituent(s);

ring D is a benzene ring which may have a substituent(s), a pyrazole ring which may have a substituent(s) or a pyrrole ring which may have a substituent(s);



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is methylene which may be substituted, ethylene which may be substituted, propylene which may be substituted, or ethenylene which may be substituted; and

Z is -COOH; -CONHSO<sub>2</sub>R<sup>1</sup>, in which R<sup>1</sup> is methyl-C1-8 alkyl which may be substituted, a benzene ring which may have a substituent(s).

- 11. (Original) The compound according to claim 1, which is selected from the group consisting of:
- (1) {1-[(2S,3S)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-(3,5-dimethoxy-4-methylphenyl)-3-hydroxylpropyl]-1H-pyrrol-3-yl}acetic acid,
- (2) (1-{(2S)-2-[(S)-(3,5-dimethoxy-4-methylphenyl)(hydroxy)methyl]-5-thien-3ylpentyl}-1H-pyrrol-3-yl)acetic acid,
- (3) {1-[(2S,3S)-2-(1,3-benzodioxol-2-ylmethyl)-3-(3,5-dimethoxy-4-methylphenyl)-3-hydroxylpropyl]-1H-pyrrol-3-yl}acetic acid,
- (4) {1-[(2S,3S)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-hydroxy-3-(3,4,5-trimethoxyphenyl)propyl]-1H-pyrrol-3-yl}acetic acid,
- (5) {1-[(2S,3S)-3-(4-acetyl-3,5-dimethoxyphenyl)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-hydroxylpropyl]-1H-pyrrol-3-yl}acetic acid,
- (6) {1-[(2S,3S)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-(4-ethyl-3,5-dimethoxyphenyl)-3-hydroxylpropyl]-1H-pyrrol-3-yl}acetic acid,
- (7) 3-{1-[(2S,3S)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-(3,5-dimethoxy-4-methylphenyl)-3-hydroxylpropyl]-1H-pyrrol-3-yl}propanoic acid,

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 $(8) \qquad 3-\{1-[(2S,3S)\cdot2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-hydroxy-3-(3,4,5-trimethoxyphenyl)propyl]-1H-pyrrol-3-yl\}propanoic acid,$ 

- (9) 3-{1-[(2S,3S)-3-(4-acetyl-3,5-dimethoxyphenyl)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-hydroxylpropyl]-1H-pyrrol-3-yl}propanoic acid,
- (10) 3-{1-[(2S,3S)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-(4-ethyl-3,5-dimethoxyphenyl)-3-hydroxylpropyl]-1H-pyrrol-3-yl}propanoic acid,
- (11) 2-{1-[(2S,3S)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-(3,5-dimethoxy-4-methylphenyl)-3-hydroxylpropyl]-1H-pyrrol-3-yl}-N-(methylsulfonyl)acetamide,
- (12) [1-[(2S,3S)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-(3,5-dimethoxy-4-methylphenyl)-3-hydroxylpropyl]-4-(methoxylcarbonyl)-1H-pyrrol-3-yl]acetic acid,
- (13) N-(3-{1-[(2S,3S)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-(3,5-dimethoxy-4-methylphenyl)-3-hydroxylpropyl]-1H-pyrrol-3-yl}propanoyl)-2-methylbenzenesulfonamide,
- (14) (2E)-3-{1-[(2S,3S)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-(3,5-dimethoxy-4-methylphenyl)-3-hydroxylpropyl]-1H-pyrrol-3-yl}acrylic acid,
- (15) 2-{1-[(2S,3S)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-(3,5-dimethoxy-4-methylphenyl)-3-hydroxylpropyl]-1H-pyrol-3-yl}-2-methylphenyl)-3-hydroxylpropyl]-1H-pyrol-3-yl}-2-methylphenyl)-3-hydroxylpropyl
- (16) (2E)-3-{1-[(2S,3S)-2-(2,3-dihydro-1H-inden-2-ylmethyl)-3-(3,5-dimethoxy-4-methylphenyl)-3-hydroxylpropyl]-1H-pyrrol-3-yl}-2-methylphenyl cacid.
- (Previously presented) A pharmaceutical composition comprising the compound of formula (I) according to claim 1, or a salt thereof, or a solvate thereof.

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13. (Original) The pharmaceutical composition according to claim 12, which is an LPA

receptor antagonist, wherein the LPA receptor is EDG-2.

14. (Canceled)

15. (Original) The pharmaceutical composition according to claim 12, which is an

agent for prevention and/or treatment for urinary system disease.

16. (Currently Amended) A method for treatment of urinary system disease selected

from the group of prostatic hypertrophy, neurogenic bladder dysfunction disease, dysuria,

pollakiuria, night urination and urodynia, which comprises administering to a mammal an

effective amount of the compound of formula (I) according to claim 1, or a salt thereof, or a

solvate thereof.

17. (Canceled)

18. (Currently Amended) A pharmaceutical composition comprising a combination of

the compound of formula (I) according to claim 1, or a salt thereof, a solvate thereof with at least

one agent selected from an LPA receptor antagonist, an al blocking agent, an anticholinergic

agent, a 5\alpha-reductase inhibitor and an anti-androgenic agent.

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